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# AN ALGORITHM FOR COMPUTING REDUCING SUBSPACES BY BLOCK DIAGONALIZATION

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## Abstract

This paper describes an algorithm for reducing a real matrix A to block diagonal form by a real similarity transformation. The columns of the transformation corresponding to a block span a reducing subspace of A, and the block is the representation of A in that subspace with respect to the basis. The algorithm attempts to control the condition of the transformation matrices, so that the reducing subspaces are well conditioned and the basis vectors are numerically independent.

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# AN ALGORITHM FOR COMPUTING REDUCING SUBSPACES BY BLOCK DIAGONALIZATION

Connice A. Bavely
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## 1. Introduction

The purpose of this report is to describe an algorithm for reducing a real matrix A of order n to a block diagonal form by a real similarity transformation. Specifically, the algorithm attempts to compute a real nonsingular matrix X such that  $X^{-1}AX$  has the form

(1.1) 
$$X^{-1}AX = B = diag(B_1, B_2, ..., B_s)$$
,

where each matrix  $B_i$  is square of order  $n_i$ .

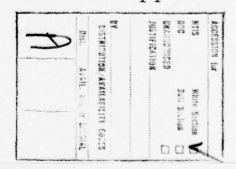
A decomposition such as (1.1) has many applications. When the blocks B, are small, powers of A can be economically calculated in the form

$$A^{k} = X \operatorname{diag}(B_{1}^{k}, B_{2}^{k}, \dots, B_{S}^{k})X^{-1}$$
,

and this fact can be used to simplify the computation of functions of A defined by power series (e.g. see [7]). If X is partitioned in the form

$$X = (X_1, X_2, ..., X_s)$$
,

where each  $X_i$  has  $n_i$  columns, then the columns of  $X_i$  form a basis for a reducing subspace of A, and  $B_i$  is the representation of A with respect to that basis. The associated spectral projector is given by  $X_i X_i^{(-1)}$ , where



 $X_i^{(-1)}$  is formed from the corresponding rows of  $X^{-1}$  (for definitions and applications see [4]).

There are theoretical and practical limitations on how small the blocks in (1.1) can be. Theoretically, they can be no smaller than the blocks in the Jordan canonical form of A. Practically, they may have to be larger. The numerical problems associated with decompositions such as (1.1) have been examined in detail in [3]. Here we give only a brief summary.

The principal difficulty is that the Jordan form of a matrix need not be numerically well determined; very small perturbations in the matrix may cause blocks to split or coalesce. Any attempt to separate two such "nearby" blocks will result in a transformation matrix X whose columns are nearly linearly dependent, or equivalently X will be ill-conditioned in the sense that the product  $\|X\|\|X^{-1}\|$  is large (here  $\|\cdot\|$  denotes a suitable matrix norm). In this case, it will be impossible to form  $X^{-1}$  or solve linear systems involving X accurately [10,14]. The phenomenon is loosely associated with close eigenvalues; but there are matrices with equal eigenvalues, e.g. symmetric matrices, that can be split completely into 1 × 1 blocks, and there are matrices with well separated eigenvalues that cannot be split except by very ill-conditioned transformations.

Our algorithm attempts to avoid these difficulties by working only with well-conditioned transformations. If a group of eigenvalues cannot be split off into a block by a transformation whose condition observes a tolerance provided by the user, the block is enlarged until a well-conditioned reducing transformation can be found. In principle this does not insure that the final transformation will be well-conditioned, since it is formed as the product of a number of reducing transformations; however, we have found that when a matrix possesses a well-conditioned decomposition of the form (1.1), our algorithm generally finds it. And the exceptions have not so much to do with the ill-conditioning of X as with the failure of the algorithm to split the matrix completely owing to the comingling of degenerate eigenvalues with well-conditioned ones.

A good deal of work has been done on the numerically stable simplification of matrices by similarity transformations [5,6,8,12,13], most of which has been summarized in [3]. For the most part, these algorithms attempt to go farther than ours in reducing the matrix, however at considerable cost in complexity and computation. The virtues of the algorithm proposed here are its simplicity and economy. When it is required to reduce a matrix beyond what is done by our algorithm, the other techniques can be applied to the blocks produced by our algorithm. The algorithm also has the advantage that it works entirely with real matrices by the device of grouping pairs of complex conjugate eigenvalues in the same block.

In the next section of this paper the algorithm is described. In Section 3 some numerical examples are given. Programming details and a listing are given in an appendix.

## 2. The algorithm.

The first part of the algorithm uses orthogonal transformations to reduce the matrix A to quasi-triangular form, that is to a block upper-triangular form in which the diagonal blocks are of order at most two. The blocks of order one contain real eigenvalues of A and the blocks of order two contain complex conjugate pairs of eigenvalues. The ordering of the blocks is arbitrary, and the order can be changed by applying appropriate orthogonal transformations. Since this reduction of A can be effected by standard techniques [9,11], we may assume that A is already in quasi-triangular form.

The subsequent block diagonalization is accomplished as follows. The matrix A is partitioned in the form

$$A = \begin{pmatrix} A_{11} & A_{12} \\ 0 & A_{22} \end{pmatrix},$$

where initially  $A_{11}$  is  $1 \times 1$  or  $2 \times 2$  depending on the dimension of the leading diagonal block of A. An attempt is then made to find a similarity transformation X such that

$$X^{-1}AX = \begin{pmatrix} A_{11} & 0 \\ 0 & A_{22} \end{pmatrix}.$$

If such a transformation can be found and if it is not too ill-conditioned, the reduction proceeds with the submatrix  $A_{22}$ . If not, a suitable  $1 \times 1$  or  $2 \times 2$  diagonal block from  $A_{22}$  is located and moved by means of orthogonal transformations to the leading position of  $A_{22}$ . The block is then

adjoined to A<sub>11</sub> by increasing the order of A<sub>11</sub> by one or two, as is appropriate, and another attempt is made to find a reducing matrix X.

The implementation of such an algorithm requires the answers to two questions.

- 1. How may the transformation X be computed?
- 2. In the event of failure, which block of A<sub>22</sub> is to be incorporated into A<sub>11</sub>?

We shall now answer these questions.

We seek the transformation X in the form

$$X = \begin{pmatrix} I & P \\ 0 & I \end{pmatrix},$$

where the identity matrices are of the same orders as  $A_{11}$  and  $A_{22}$ . The inverse of X is easily seen to be

$$(2.2) X^{-1} = \begin{pmatrix} I & -P \\ 0 & I \end{pmatrix}.$$

Hence

$$X^{-1}AX = \begin{pmatrix} A_{11} & A_{11}^{P-PA} & A_{22}^{+A} & A_{22} \end{pmatrix},$$

and the problem of determining X becomes that of solving the equation

$$A_{11}^{P} - PA_{22} = -A_{12}.$$

Because  $A_{11}$  and  $A_{22}$  are quasi-triangular, this equation can be solved by a back-substitution algorithm of Bartels and Stewart [1], provided the

eigenvalues of  $A_{11}$  and  $A_{22}$  are disjoint.

From (2.1) and (2.2) it follows that X will be ill-conditioned whenever P is large. As each element of P is generated, it is tested to see if its magnitude exceeds a bound provided by the user. If it does, the attempt to compute X is abandoned and a new, larger block A<sub>11</sub> is formed. If no element of P exceeds the bound, the matrix X is accepted and the matrix A is deflated as described above. The transformation X is postmultiplied into a matrix that accumulates all the transformations made on the matrix A.

The process for selecting a  $1 \times 1$  or  $2 \times 2$  block of  $A_{22}$  to incorporate into  $A_{11}$  goes as follows. We compute the mean of those eigenvalues of  $A_{11}$  having nonnegative imaginary part. A block is chosen from  $A_{22}$  whose eigenvalue with nonnegative imaginary part is nearest this mean. This block is moved, as described above, by orthogonal transformations to the leading position in  $A_{22}$ , where it is incorporated into  $A_{11}$ . The program HQR3 [11], which can be used to obtain the initial quasitriangular form, has a subroutine which will compute these orthogonal transformation. The transformations are of course postmultiplied into the accumulating matrix.

We summarize our algorithm in the following informal code. Further details can be found in the appendix to this paper, where a FORTRAN subroutine implementing this code is given. The code takes as input an array A of order N containing the matrix A and an array X in which the transformations are accumulated. In addition the user must provide a

tolerance to bound the size of the elements of the deflating transformations. The integers L11 and L22 point to the beginnings of the current blocks  $A_{11}$  and  $A_{22}$  in the array A. The informal code should be self-explanatory. Comments are delineated by the symbol #.

```
reduce A to quasitriangular form, accumulating the
              transformations in X [10];
            L11 = 1;
            loop # until the matrix is diagonalized #
3
3.1
              if L11 > N then leave 3 fi;
              L22 = L11;
3.2
              loop # until a block has been deflated #
3.3
                 if L22 = L11 then # use the first 1 × 1 or 2 × 2 block #
3.3.1
                   M = order of the block at L11;
3.3.1t.1:
                   L22 = L22 + M;
3.3.1t.2:
                 else # augment A_{11} with a 1 × 1 or 2 × 2 block from A_{22} # compute the mean of the eigenvalues of A_{11} with
3.3.1
3.3.le.1:
                     nonnegative imaginary parts;
                   find the M \times M # M = 1 or 2 # block of A_{22} whose eigen-
3.3.1e.2:
                     value with nonnegative imaginary part is nearest the
3.3.1e.3:
                   move the block to the leading position of A22 accumulating
                     the transformations in X;
                   L22 = L22 + M \# which incorporates the block in A_{11} \#;
3.3.le.4:
3.3.1
                 fi;
                if L22 > N then leave 3.3 fi;
attempt to split off A, [1];
if the attempt was successful then leave 3.3 fi;
3.3.2
3.3.3
3.3.4
                 restore A<sub>12</sub>;
3.3.5
3.3
              end loop;
3.4
               if L22 ≤ N then accumulate the deflating transformation in X
3.5
               scale columns L11 through L22-1 of X so that they have
                 2-norm unity, adjusting A accordingly;
3.6
              L11 = L22;
            end loop;
```

Several comments should be made about the algorithm. First, it uses only real arithmetic, even when A has complex eigenvalues. Second, the algorithm cannot guarantee that the final transformation is well conditioned, since the bound on the elements of P restricts the condition of

only the individual transformations comprising the final one. Nonetheless, we have found little tendency toward excessive growth in the condition of the transformation. Third, no attempt is made to segregate nearly equal eigenvalues initially into clusters; whether an eigenvalue splits off or not depends entirely on its determining a suitably small matrix P. This is important, since it means that the algorithm can compute well conditioned eigenvectors for multiple eigenvalues (a little help is required from rounding error; see Section 3).

The strategy for determining what eigenvalues to add to the current group has the defect that it can mix well-conditioned and ill-conditioned eigenvalues that are nearly equal, thus missing the possibility of a more complete reduction. This is not a serious problem when the blocks are small. However, if a finer resolution of the structure of the matrix is required, the techniques referenced in Section 1 may be applied to the blocks produced by our algorithm. In fact our algorithm can be regarded as a preprocessing step to reduce the problem to a size where it can be attacked by more sophisticated, but more expensive methods.

We note that the output of our algorithm may depend on the user supplied tolerance for the elements of P. In general the larger the tolerance, the smaller the blocks but the more ill conditioned the transformation. This tradeoff is an inevitable consequence of the poor determination of the structure of a matrix in the presence of errors, and we know of no algorithm that relieves the user of the necessity of making a decision of this kind.

So far as storage is concerned, the algorithm requires  $2n^2$  locations to contain the matrices A and X and a number of arrays of order n. This is the same as the storages required by algorithms that compute the eigenvectors of a general matrix.

Excluding the initial reduction of A to quasitriangular form, the bulk of the computations in the algorithm occur at statement 3.3.3, where an attempt is made to solve the equation (2.3), and at statement 3.4, where the transformation is accumulated. The multiplication count for the algorithm for solving (2.3) is of order  $(\ell^2 m + m^2 \ell)/2$ , where  $\ell$  is the size of  $\ell^2 m$  and m is the size of  $\ell^2 m$ . The cost of accumulating this transformation in X is of order  $\ell^2 m$ . Thus, at one extreme, if all the eigenvalues of A are real and they all can be deflated, the cost in multiplications will be of order  $\ell^2 m$ , which compares favorably with algorithms for computing the eigenvectors of A from its quasitriangular form. On the other hand, if the eigenvalues of A are real and none of them can be deflated, the algorithm will require on the order of  $\ell^4 m$ 

Although an n<sup>4</sup> operation count is disturbing, there are several mitigating factors. First, we do not expect the algorithm to find many applications to matrices that cannot be substantially reduced by it, since the object of using it is to save on subsequent calculations with the matrix. Second, the count assumes that the algorithm for solving (2.3) must be executed to completion before it is found that P is unacceptably large. This is not likely; in fact because the algorithm [11] for reducing

A to quasitriangular form arranges the eigenvalues in decreasing order of magnitude, it is rather unlikely. Finally the order constant 1/12 is is modest; for n less than 60 the bound is less than  $5n^3$ .

## 3. Numerical results.

In this section we summarize the results of some numerical experiments. Two of the tests were performed with a class of test matrices generated as follows. Let J be a given matrix whose structure is known (e.g. J could be in Jordan canonical form). Let

$$H_1 = I - \frac{2}{n} e e^T,$$

where  $e = (1, 1, ..., 1)^{T}$ , and

$$H_2 = I - \frac{2}{n} ff^T,$$

where  $f = (1,-1,...,(-1)^{n-1})^T$ . Note that  $H_1$  and  $H_2$  are symmetric and orthogonal. Let

$$S = diag(1, \sigma, \sigma^2, \dots, \sigma^{n-1})$$

where  $\sigma$  is a given parameter. Then we take A in the form

(3.1) 
$$A = (H_2SH_1)J(H_2SH_1)^{-1} = (H_2SH_1)J(H_1S^{-1}H_2).$$

The matrix A, which is cheap to compute, has the same structure as J. The transformation  $H_2SH_1$  can be made as ill-conditioned as desired by varying  $\sigma$ .

In describing the results of the tests we report two numbers. The first is the maximum  $\rho$  of the scaled residuals

$$\rho_{i} = \frac{\|AX_{i}^{-XB_{i}}\|_{\infty}}{\|A\|_{\infty}\|X_{i}\|_{\infty}},$$

where  $X_i$  is composed of the columns of X corresponding to the block  $B_i$ . Second we report  $\|X^{-1}\|_{\infty}$ . Together these numbers give an idea of how stably the reduction proceeded; for if

$$R = AX - XB$$

then, with  $E = RX^{-1}$ , we have that

$$(A+E)X - XB = 0 ;$$

that is B is exactly similar to A + E. The relative error that this perturbation represents in A is

$$\frac{\|E\|}{\|A\|} \leq \frac{\|R\|\|X^{-1}\|}{\|A\|} .$$

Since  $\|X\| \cong 1$ , the relative error will be of the order  $\|\rho\|X^{-1}\|$  .

The first test case illustrates the ability of the algorithm to split apart nearly equal eigenvalues with independent eigenvectors. We took

$$J = diag[1,1-\epsilon,1+\epsilon, \begin{pmatrix} 1 & -1 \\ 1 & 1 \end{pmatrix}, .3, .4, .5, .6, .7]$$
.

The algorithm was applied for various values of  $\epsilon$ ,  $\sigma$ , and RMAX, the bound

on the size of the deflating transformations. The results are summarized in Table 1, in which the eigenvalues of A are numbered as follows:

Complex eigenvalues are indicated by a circumflex.

Provided RMAX was large enough to allow a sufficiently ill conditioned transformation, all the cases were split completely. The condition of the transformation was in no case much greater than the condition of the scrambling transformation in (3.1). It is of interest to note that the algorithm was successful when  $\varepsilon = 0$ , that is when A has three equal eigenvalues. Mathematically, the algorithm for solving (2.3) breaks down when  $A_{11}$  and  $A_{22}$  have common eigenvalues; however, the experiments indicate that if the equal eigenvalues have independent vectors, rounding error will perturb them enough for the algorithm to work.

The second example shows the failure of the algorithm's strategy for selecting the next eigenvalue to be adjoined to  $A_{11}$ . Here J has the form

$$J = \operatorname{diag} \left[ 1, \begin{pmatrix} 1 & 1 & 0 \\ 0 & 1 & 1 \\ 0 & 0 & 1 \end{pmatrix}, .3, .4, .5, .6, .7, .8 \right]$$

and  $\sigma$  was taken to be one. Of the four eigenvalues at unity, one is perfectly conditioned, and the other three, which belong to a single Jordan block, are very ill-conditioned. To six figures the computed eigenvalues were

1.00073 + 0.001303i 1.00073 - 0.001303i 1.00000 0.99853

.80000

The pair of complex eigenvalues could not be deflated, since they were coupled to the third member of the block. But this member would not be adjoined without first adjoining the well conditioned eigenvalue at unity. Consequently, the algorithm produced a single block of order four, rather than two blocks of orders one and three. This block of order four could be reduced further by the more sophisticated techniques described in the references.

The third test case is the Frank matrix of order 12 which has appeared frequently in the literature [2,3]. The smaller eigenvalues of this matrix are extremely ill conditioned. The results of test runs on this matrix are summarized below.

RMAX	BLOCK STRUCTURE	ρ	$\ \mathbf{x}^{-1}\ _{\infty}$
10	1,2,3,4,5,(6,7,8,9,10,11,12)	$2.9 \times 10^{-7}$	93.6
50	1,2,3,4,5,6,(7,8,9,10,11,12)	$2.9 \times 10^{-7}$	2920.2
100	1,2,3,4,5,6,(7,8,9,10,11,12)	$2.9 \times 10^{-7}$	2920.2
1000	1,2,3,4,5,6,(7,8,9,10,11,12)	$2.9 \times 10^{-7}$	2920.2

The algorithm performed much as expected, separating the larger eigenvalues

and grouping the smaller eigenvalues together. This grouping is consistent with the precision of the computation.

The final test case is included because the failure of our algorithm to decompose it reveals the shaky foundations of a fairly common numerical practice. Specifically we generated the companion matrix of the polynomial given in [14, p. 74]. Since the zeros of this polynomial are not very ill conditioned, we were surprised when the algorithm failed to split off so much as one. Some further computations revealed that the matrix of left eigenvectors (the inverse Vandermode of the zeros) had rows of order  $10^6 - 10^8$ . This explains the failure to reduce the matrix. More important, though, it shows that the eigenvalues of the companion matrix are much more ill conditioned than the zeros of the polynomial and suggests that the practice of using eigenvalue routines to find zeros of polynomials can result in an unnecessary loss in accuracy.

σ	ε	RMAX	BLOCK STRUCTURE	ρ   X <sup>-1</sup>    <sub>∞</sub>
1.0	10 <sup>-1</sup>	10.	(1,2),3,4,5,6,7,8,9,10	$2.1 \times 10^{-7}$ 7.7
	10 <sup>-5</sup>	10.	(1,2),3,4,5,6,7,8,9,10	$1.4 \times 10^{-7}$ 6.7
	0.0	10.	(1,2),3,4,5,6,7,8,9,10	$1.1 \times 10^{-7}$ 8.0
10.0	10 <sup>-1</sup>	10.	(1,2),3,4,5,6,7,8,9,10	$1.5 \times 10^{-7}$ 25.6
	10 <sup>-5</sup>	10.	(1,2),3,4,5,6,7,8,9,10	$1.8 \times 10^{-7}$ 25.4
	0.0	10.	(1,2),3,4,5,6,7,8,9,10	$1.6 \times 10^{-7}$ 24.2
100.0	10 <sup>-1</sup>	10.	(1,2),3,4,5,6,7,8,9,10	$1.1 \times 10^{-7}$ 187.9
ody To n	10 <sup>-5</sup>	10.	(1,2),3,4,5,6,7,8,9,10	$1.3 \times 10^{-7}$ 196.6
	0.0	10.	(1,2),3,4,5,6,7,8,9,10	$1.2 \times 10^{-7}$ 157.9
1000.0	10 <sup>-1</sup>	10.	(1,2,4,3,5),(6,7,8,9),10	$1.1 \times 10^{-7}$ 326.6
		100.	(1,2),3,4,5,6,7,8,9,10	$1.0 \times 10^{-8}$ 1537.9
	10 <sup>-5</sup>	10.	(1,2,4,3,5),(6,7,8,9),10	$7.8 \times 10^{-8}$ 333.8
		100.	(1,2),3,4,5,6,7,8,9,10	$7.8 \times 10^{-8}$ 1559.1
	0.0	10.	(1,2,3),4,5,(6,7,8,9),10	$8.2 \times 10^{-8}$ 394.1
		100.	(1,2),3,4,5,6,7,8,9,10	$8.2 \times 10^{-8}$ 1355.7

Table 4.1

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# Appendix

#### PROGRAMMING DETAILS AND PROGRAM LISTING

## Al. Usage.

The calling sequence for BDIAG is

CALL BDIAG(A,LDA,N,EPSHQR,RMAX,ER,EI,TYPE,BS,X,LDX,FAIL).

The parameters in the calling sequence are (starred parameters are altered by the subroutine)

*A(LDA,N)	an array that initially contains the $N \times N$ matrix
	to be reduced. On return A contains the reduced
	block diagonal matrix.

LDA the leading dimension of A.

N the order of the matrices A and X.

EPSHQR a real variable containing a convergence criterion for the subroutines HQR3 and EXCHNG [11].

RMAX a real variable containing a bound on the absolute values of the elements in the reducing matrices.

\*ER(N) a real array containing the real parts of the eigenvalues of A.

\*EI(N) a real array containing the imaginary parts of the eigenvalues.

\*TYPE(N) an integer array whose i-th entry is 0 if the i-th eigenvalue is real

1 if the i-th eigenvalue is complex with positive imaginary part

2 if the i-th eigenvalue is complex with negative imaginary part

-1 if the i-th eigenvalue could not be computed

\*BS(N) a singly subscripted array that contains information on the block structure of the matrix returned by the program. If there is a block of order K at A(L,L), then BS(L),BS(L+1),...BS(L+K-1) contain the integers K,-(K-1),...,-1. Thus a positive entry of K indicates the start of a block of order K.

\*X(LDX,N) an array into which the reducing transformations are accumulated.

LDX the leading dimension of X.

FAIL a logical variable which is false on a normal return and is true on return if an error has occurred.

The program requires the programs ORTHES [9], ORTRAN [9], HQR3 [11], EXCHNG [11], SHRSLV [1], DAD and SPLIT [11].

A suitable choice for the parameter EPSHQR is the rounding unit of the computer on which the program is being run; i.e. if one is working in a precision of about t decimal figures EPSHQR should be of order 10<sup>-t</sup>.

A2. Programming details.

In this section we shall describe some of the details of the implementation of the algorithm. Throughout this section we refer to the outline of the algorithm in Section 2.

Statement 1. This is accomplished by the EISPACK routines ORTHES and ORTRAN [9] and the QR routine HQR3 [11].

Statement 2. L11 points to the leading position of the current block A11, which is of order DA11.

Statement 3. This is the main loop of the program. It ends when L11 > N, indicating that there are no further blocks to deflate.

Statement 3.2. L22 points to the leading position of the current block A22, which is of order DA22.

Statement 3.3. In this loop  $A_{11}$  is repeatedly augmented until it can be deflated or until  $A_{22}$  is void.

Statements 3.3.1t.  $A_{11}$  is initially void. Here it is taken to be the 1 × 1 or 2 × 2 block starting at L11.

Statements 3.3.1e. This is the search for a  $1 \times 1$  or  $2 \times 2$  block described in Section 2.

Statement 3.3.1e.3. The subroutine EXCHNG [11] is used repeatedly to move the block just located to the beginning of  $A_{22}$ . After each exchange of a complex block SPLIT [11] is called to recompute its eigenvalues and to see if, owing to rounding error, it can be split into a pair of real eigenvalues.

Statement 3.3.2. Because A22 is void, A11 is effectively deflated.

Statement 3.3.3. The matrix  $A_{12}$  is saved below the lower subdiagonal of A, in case the attempt to deflate  $A_{11}$  is unsuccessful. Since the routine SHRSLV [1], which computes the deflating transformation, requires  $A_{11}$  to be lower Hessenberg, the subroutine DAD is called to transform  $A_{11}$  and  $A_{12}$ . SHRSLV has been modified to return with a signal if some element of the deflating transformation exceeds the bound RMAX. Otherwise the matrix P that determines the transformation overwrites  $A_{12}$ . DAD is once again called to restore  $A_{11}$  to its original form.

Statement 3.3.5. The submatrix  $A_{22}$ , which was overwritten by SHRSLV, must be restored before another attempt to deflate is made.

Statement 3.4. Only if  $A_{22}$  is not void was a deflating transformation generated.

Statement 3.6. Set L11 to point to the next submatrix before continuing.

# Program listing.

SUBROUTINE BDIAG (A.LDA.N.EPSHOR.RMAX.ER,EI,TYPF,BS,X,LDX,FAIL)

INTEGER LDA, LDX, N, TYPE(N), BS(N)
REAL A(LDA,N), ER(N), EI(N), X(LDX,N), EPSHOR, RMAX
LOGICAL FAIL

HDIAG REDUCES A MATRIX A TO BLOCK DIAGONAL FORM BY FIRST REDUCING IT TO QUASI-TRIANGULAR FORM BY HOP3 AND THEN BY SOLVING THE MATRIX EQUATION -A11\*P+P\*A22=A12 TO INTRODUCE ZEROS ABOVE THE DIAGONAL. THE PARAMETERS IN THE CALLING SEQUENCE ARE (STARRED PARAMETERS ARE ALTERED BY THE SUBROUTINE):

AN ARRAY THAT INITIALLY CONTAINS THE N X N MATRIX TO BE REDUCED. ON RETURN, A CONTAINS THE REDUCED BLOCK DIAGONAL MATRIX. \* A

THE LEADING DIMENSION OF ARRAY A. LDA

THE ORDER OF THE MATRICES A AND X.

THE CONVERGENCE CRITERION FOR SUBROUTINE HOR3. EPSHOR

THE MAXIMUM SIZE ALLOWED FOR ANY ELEMENT OF THE TRANSFORMATIONS. RMAX

A SINGLY SUBSCRIPTED REAL ARRAY CONTAINING THE REAL PARTS OF THE EIGENVALUES. \*FR

A SINGLY SUBSCRIPTED REAL ARRAY CONTAING THE IMAGINAPY PARTS OF THE EIGENVALUES. \*LI

SINGLY SUBSCRIPTED INTEGER ARRAY WHOSE I-TH ENTRY IS IF THE I-TH EIGENVALUE IS REAL IF THE I-TH EIGENVALUE IS COMPLEX WITH POSITIVE \*TYPE

IMAGINARY PART

IF THE I-TH EIGENVALUE IS COMPLEX WITH POSITIVE

IMAGINARY PART

IMAGINARY PART

THE I-TH EIGENVALUF COULD NOT BE COMPUTED

A SINGLY SUBSCRIPTED INTEGER ARRAY THAT CONTAINS BLOCK STRUCTURE INFORMATION. IF THERE IS A BLOCK OF ORDER K STARTING AT A(L,L) IN THE OUTPUT MATRIX A, THEN BS(L) CONTAINS THE POSITIVE INTEGER K, BS(L+1) CONTAINS \*B5 THUS A POSITIVE INTEGER IN THE L-TH FNTRY OF BS INDICATES A NEW BLOCK OF ORDER BS(L) STARTING AT A(L,L).

AN ARRAY INTO WHICH THE REDUCING TRANSFORMATIONS ARE TO BE MULTIPLIED.

THE LEADING DIMENSION OF ARRAY X. LDX

A LOGICAL VARIABLE WHICH IS FALSE ON NORMAL RETURN AND TRUE IF THERE IS ANY ERROR IN BDIAG. \*FAIL

BDIAG USES SUPROUTINES ORTHES, ORTRAN, HOR3, EXCHNG, SHRSLV, DAD, AND SPLIT.

INTEGER DA11, DA22, I, J, K, KM1, KM2, L, LFAVE, LOOP, L11, L22, L22M1, NK REAL C, CAV, D, E1, E2, RAV, SC, TEMP

FAIL = , TRUE.

CONVERT A TO UPPER HESSENHERG FORM.

CALL ORTHES (LDA.N. 1.N. A. ER)

CCC

a

\* X

```
CALL ORTRAN (LDA,N,1,N,A,ER,X)
CCC
           CUNVERT A TO QUASI-UPPER TRIANGULAR FORM BY OR METHOD.
           CALL HOR3 (A,X,N,1,N,EPSHOR,ER,EI,TYPE,LDA,LDX)
0000
           CHECK TO SEE IF HOR3 FAILED IN COMPUTING ANY EIGFNVALUE
           DU 5 I = 1.N
IF (TYPE(I).EQ.-1) GO TO 900
           CONTINUE
     5
REDUCE A TO BLOCK DIAGONAL FORM
           SEGMENT A INTO 4 MATRICES: All, A DAll X DALL BLOCK WHOSE (1,1)-ELEMENT IS AT A(L11,L11); A22, A DA22 X DA22 BLOCK WHOSE (1,1)-ELEMENT IS AT A(L22,L22); A12, A DA11 X DA22 BLOCK WHOSE (1,1)-ELEMENT IS AT A(L11,L22); AND A21, A DA22 X DA11 BLOCK = 0 WHOSE (1,1)-ELEMENT IS AT A(L22,L11).
           THIS LOOP USES L11 AS LOOP INDEX AND SPLITS OFF A BLOCK STARTING AT A(L11,L11).
           L11 = 1
ASSIGN 550 TO LOOP
ASSIGN 600 TO LEAVE
IF (L11 .GT. N) GO TO LEAVE
L22 = L11
    10
COCO
               THIS LOOP USES DA11 AS LOOP VARIABLE AND ATTEMPTS TO SPLIT OFF A BLOCK OF SIZE DA11 STARTING AT A(L11,L11)
               ASSIGN 350 TO LOOP

IF (L22 • NE • L11) GO TO 110

DA11 = TYPE(L11) + 1

L22 = L11 + DA11

L22M1 = L22 - 1

GO TO 290
   100
                   CONTINUE
   110
COOO
                       COMPUTE THE AVERAGE OF THE EIGENVALUES IN A11
                       RAV = 0.

CAV = 0.

DO 120 I = L11.L22M1

RAV = RAV + ER(I)

CAV = CAV + ARS(EI(I))

CONTINUE

RAV = RAV / FLOAT(DA11)

CAV = CAV / FLOAT(DA11)
    120
000
                       LOOP ON EIGENVALUES OF A22 TO FIND THE ONE CLOSEST TO THE AVERAGE.
                       D = (RAV-ER(L22))**2 + (CAV-EI(L22))**2
K = L22
L = L22 + TYPE(L22) + 1
                       L = L22 + TYPF(L22) + 1

ASSIGN 145 TO LOOP

IF (L.GT.N) GO TO 150

C = (RAV-ER(L))**2 + (CAV-EI(L))**2

IF (C.GE.D) GO TO 140
    130
                           CONTINUE TYPE(L) + 1
   140
                       GO TO 130
CONTINUE
   145
0000
                       LOOP TO MOVE THE EIGENVALUE JUST LOCATED INTO FIRST POSITION OF BLOCK A22.
```

```
C
                    ASSIGN 280 TO LOOP
IF (TYPE(K).NE.0) GO TO 200
                        THE BLOCK WE'RE MOVING TO ADD TO All IS A 1 X 1
                       NK = 1
CONTINUE
IF (K .EQ. L22) GO TO 280
KM1 = K - 1
   160
                        IF (TYPE(KM1).EQ.0) GO TO 190
                       WE'RE SWAPPING THE CLOSEST BLOCK WITH A 2 X 2
                       KM2 = K - 2
CALL EXCHNG (A,X,N,KM2,2,1,EPSHQR,FAIL,1DA,LDX)
IF (FAIL) GO TO 900
                       TRY TO SPLIT THIS BLOCK INTO 2 REAL EIGENVALUES
                       CALL SPLIT (A.X.N.KM1,F1,E2,LDA,LDX) IF (A(K.KM1).EQ.O.) GO TO 170
                    (M2) = 0

(KM1) = 1

(PE(K) = 2

ER(KM2) = 0.

ER(KM2) = 0.

ER(KM1) = E1

ER(KM1) = E2

ER(KM1) = E2

EX(KM1) = E2
                       BLOCK IS STILL COMPLEX.
                       COMPLEX BLOCK SPLIT INTO TWO REAL EIGENVALUES.
                       CONTINUE

TYPE(KM1) = 0

TYPE(KM2) = 0

ER(KM2) = ER(K)

ER(KM1) = E1

ER(K) = E2

EI(KM1) = 0
   170
                       FI(KMI) = 0.

K = KM2

IF (K .LE. L22) GO TO 280

GO TO 160
   180
CCCC
                       WE'RE SWAPPING THE CLOSEST BLOCK WITH A 1 X 1.
                       CONTINUE
CALL EXCHAG (4, X, N, KM1, 1, 1, EPSHOR, FAIL, LDA, LDX)
IF (FAIL) GO TO 900
TEMP = ER(K)
ER(K) = ER(KM1)
ER(KM1) = TEMP
   190
                       K = KM1
                       F (K LE. L22) GO TO 280
                       THE BLOCK WE'RE MOVING TO ADD TO All IS A 2 X 2.
                       CONTINUE
CONTINUE
   200
   210
                       IF (K .EQ. L22) GO TO 280
                        IF (TYPE(KM1) .EQ. 0) GO TO 240
                       WE'RE SWAPPING THE CLOSEST BLOCK WITH A 2 X 2 BLOCK.
                       KM2 = K - 2
                                           (A.X.N.KM2.2.2.EPSHOR.FAIL.LDA.LDX)
```

```
IF (FAIL) GO TO 900
CCC
                             TRY TO SPLIT SWAPPED BLOCK INTO TWO REALS.
                             CALL SPLIT (A,X,N,K,E1,E2,LDA,LDX)
ER(KM2) = ER(K)
ER(KM1) = ER(K+1)
                             EI(KM2) = EI(K)

EI(KM1) = EI(K+1)
                             IF (A(K+1.K) .EQ. 0.) GO TO 220
                             STILL COMPLEX BLOCK.
                             ER(K) = E1
ER(K+1) = E1
EI(K) = E2
EI(K+1) = -E2
                             GO TO 230
000
                             TWO REAL ROOTS.
                             CONTINUE

TYPE(K) = 0

TYPE(K+1) = 0

ER(K) = E1

ER(K+1) = E2

EI(K) = 0.

EI(K+1) = 0.

CONTINUE

K = KM2

IF (K.EQ.L22) GO TO 260

GO TO 210
    220
    230
CCC
                              WE'RE SWAPPING THE CLOSEST BLOCK WITH A 1 X 1.
                             CONTINUE

CALL EXCHNG (A,X,N,KM1,1,2,EPSHQR,FAIL,1DA,LDX)

IF (FAIL) GO TO 900

TYPE (KM1) = 1

TYPE(K) = 2

TYPE(K+1) = 0

ER(K+1) = ER(KM1)

FR(KM1) = ER(K)

EI(KM1) = EI(K)

EI(KM1) = EI(K)

EI(K) = EI(K+1)

EI(K+1) = 0.

GO TO 250
    240
    250
                              CONTINUE
                              K = KM1
IF (K.EQ.L22) GO TO 260
GO TO 210
CCC
                              TRY TO SPLIT RELOCATED COMPLEX BLOCK.
                              CONTINUE
CALL SPLIT (A.X.N.K.E1.E2.LDA.LDX)
IF (A(K+1.K).EQ.O.) GO TO 270
    260
000
                              STILL COMPLEX.
                             ER(K) = E1
ER(K+1) = E1
EI(K) = E2
EI(K+1) = -E
GO TO 280
                                                  -E2
CCC
                              SPLIT INTO TWO REAL EIGENVALUES.
    270
                              CONTINUE
                             TYPE(K) = 0

TYPE(K+1) =

ER(K) = E1

ER(K+1) = E2

EI(K) = 0.
                                                 0 0
```

```
C
                CONTINUE
DA11 = DA11 + NK
L22 = L11 + DA11
L22M1 = L22 - 1
CONTINUE
ASSIGN 400 TO LEAVE
IF (L22 .GT. N) GO TO LEAVE
   280
   290
CCC
                ATTEMPT TO SPLIT OFF A BLOCK OF SIZE DA11.
                DA22 = N - L22 + 1
CCC
                SAVE A12 IN ITS TRANSPOSE FORM IN BLOCK A21.
                300
                CONTINUE
00000
                CONVERT A11 TO LOWER QUASI-TRIANGULAR AND MULTIPLY IT BY -1 AND CHANGE A12 APPROPRIATELY (FOR SOLVING -A11*P+P*A22=A12).
                CALL DAD (A,LDA,L11,L22M1,L11,N,1.,0)
CALL DAD (A,LDA,L11,L22M1,L11,L22M1,-1.,1)
CCC
                SOLVE -A11*P + P*A22 = A12.
                CALL SHRSLV (A(L11,L11),A(L22,L22),A(L11,L22),
DA11,DA22,LDA,LDA,LDA,RMAX,FAIL)
ASSIGN 400 TO LEAVE
        1
                IF (.NOT. FAIL) GO TO LEAVE
CCC
                CHANGE All BACK TO UPPER QUASI-TRIANGULAR.
                CALL DAD (A.LDA.L11.L22M1.L11.L22M1.1.1)
CALL DAD (A.LDA.L11.L22M1.L11.L22M1.-1..0)
000000
                WAS UNABLE TO SOLVE FOR P - TRY AGAIN
                MOVE SAVED A12 BACK INTO ITS CORRECT POSITION.
                DO 310 J=L11.L22M1

DO 310 I = L22.N

A(J,I) = A(I,J)

A(I,J) = 0.

CONTINUE
   310
             GO TO 100
CONTINUE
   350
CCC
             CHANGE SOLUTION TO P TO PROPER FORM.
             IF (L22 .GT. N) GO TO 440

CALL DAD (A,LDA,L11,L22M1,L11,N,1.,0)

CALL DAD (A,LDA,L11,L22M1,L11,L22M1,-1.,1)
00000
                MULTIPLY TRANSFORMATION INTO X.
                ONLY COLUMNS L22 THRU N ARF AFFECTED.
                DO 410 J = L22.N

DO 410 I = 1.N

DO 410 K = L11.L22M1

X(I,J) = X(I,J) + X(I,K) * A(K,J)
   410
                ZERO OUT A12 FOR EASE IN HANDLING.
                DO 420 J = L22,N

DO 420 I = L11,L22M1

A(I,J) = 0.

CONTINUE
   420
```

```
C
CC
                             ZERO OUT TRIANGULAR BLOCK BELOW DIAGONAL .
                            DO 430 J = L11; 22M1
DO 430 I = L22,N
A(1,J) = 0.
CONTINUE
     430
COCO
                       SCALE THOSE COLUMNS OF X THAT WON'T BE ALTERED AGAIN TO UNITY. CHANGE All APPROPRIATELY.
                      CONTINUE

DO 500 J = L11.L22M1

SC = 0.

DO 450 I = 1.N

SC = SC + (X(I,J))**2

CONTINUE

SC = SQRT(SC)

DO 460 I = 1.N

X(I,J) = X(I,J) / SC

CONTINUE

DO 470 I = L11.L22M1

A(I,J) = A(I,J) / SC

CONTINUE

DO 480 I = L11.L22M1

A(J,I) = A(J,I) * SC

CONTINUE

CONTINUE

CONTINUE

CONTINUE

CONTINUE

CONTINUE

CONTINUE

CONTINUE
      440
      450
      460
      470
      480
500
CCC
                        STORE BLOCK SIZE IN ARRAY BS.
                       BS(L11) = DA11

J = DA11 - 1

IF (J .EQ. 0) GO TO 520

DO 510 I = 1,J

BS(L11+I) = -(DA11-I)

CONTINUE
      510
520
     520 CONTINUE
L11 = L22
550 GO TO 10
600 CONTINUE
FAIL = FALSE.
RETURN
                  ERROR RETURN.
      900 CONTINUE
FAIL = •TRUE•
RETURN
END
```

```
00000000000000
                                                             IF ISW = 0, SURROUTINE DAD
COMPUTES D*A WHERE D IS THE MATRIX WITH ONES DOWN
THE MINOR DIAGONAL AND A IS THE INPUT MATRIX.

IF ISW = 1, IT COMPUTES THE PRODUCT A*D.

IT COMPUTES THIS PRODUCT FOR ROWS I1 THRU 12 AND COL-
UMNS J1 THRU J2. IT ALSO MULTIPLIES EACH ELEMENT OF
THE PRODUCT WITH THE CONSTANT R. NA IS THE FIRST
DIMENSION OF THE MATRIX A. THE PRODUCT
OVERWRITES THE SPECIFIED LOCATIONS OF MATRIX A.
                                   REAL A(NA.1), R
IF (ISW .EG. 1) GO TO 200
IF (II .EG. 12) GO TO 150
           NRD2 = IFIX((I2 - I1 + 1) / 2

DO 100 J = J1,J2

DO 50 IP1 = 1,NRD2

I = IP1 - 1

TEMP = A(I1+I,J)

A(I1+I,J) = A(I2-I,J) * R

A(I2-I,J) = TEMP * R

50 CONTINUE

IF (MOD(I2-I1,2) .EQ. 1) RETURN

I = II + NRD2

DO 110 J=J1,J2

A(I,J) = A(I,J) * R

110 CONTINUE

RETURN

150 CONTINUE

DO 160 J = J1,J2

A(I1,J) = A(I1,J) * R

160 CONTINUE

RETURN
 C
 000000
                                                         COMPUTES THE PRODUCT AD WHFRE D IS AS AROVE.
           200 CONTINUE

IF (J1 .E0. J2) GO TO 350

NCD2 = IFIX((J2 - J1 + 1) / 2

DO 300 JP1 = 1.NCD2

J = JP1 - 1

UO 250 I = II.I2

TEMP = A(I.J1+J)

A(I.J1+J) = A(I.J2-J) * R

250 CONTINUE

300 CONTINUE

IF ( MOD(J2-J1.2) .E0. 1) RETURN

J = J1 + NCD2

DO 310 I=I1.I2

A(I.J) = A(I.J) * R

310 CONTINUE

RETURN

350 CONTINUE
             350 CONTINUE

UO 360 I = I1.12

A(I.J1) = A(I.J1) * R

360 CONTINUE

RETURN
                                    END
```

SUBROUTINE DAD (A, NA, II, I2, J1, J2, R, ISW)

```
SUBROUTINE SHRSLV (A.B.C.M.N.NA.NB.NC.RMAX.FAIL)
SHRSLV IS A FORTRAN IV SUBROUTINF TO SOLVE THE REAL MATRIX EQUATION AX + XB = C, WH RE A IS IN LOWER REAL SCHUR FORM AND B IS IN UPPER REAL SCHUR FORM. SHRSLV USES THE AUX-ILLIARY SUBROUTINE SYSSLV, WHICH IT COMMUNICATES WITH THROUGH THE COMMON BLOCK SLVBLK. THE PARAMETERS IN THE CALLING SEQUENCE ARE
                        A DOUBLY SUBSCRIPTED ARRAY CONTAING THE MATRIX A IN LOWER SCHUR FORM
                        A DOUBLY SUBSCRIPTED ARRAY CONTAINING THE MATRIX B IN UPPER REAL SCHUR FORM
                        A DOUBLY SUBSCRIPTED ARRAY CONTAINING THE MATRIX C.
        C
                        THE ORDER OF THE MATRIX A
                        THE ORDER OF THE MATRIX B
        N
                        THE FIRST DIMENSION OF THE ARRAY A
        NA
        NB
                        THE FIRST DIMENSION OF THE ARRAY B
        NC
                        THE FIRST DIMENSION OF THE ARRAY C
        RMAX
                        MAXIMUM ALLOWED SIZE OF ANY ELEMENT OF THE TRANSFORMATION
                      INDICATES IF SHRSLV FAILED
        FAIL
             INTEGER M.N.NA.NB.NC.K.KM1.DK.KK.L.LM1.DL.LL.I.IB.J.JA.NSYS
REAL A(NA.1), B(NB.1), C(NC.1), T. P
LUGICAL FAIL, SING
COMMON/SLVBLK/ T(5.5),P(5),NSYS.SING
FAIL = .TRUE.
            L = 1
LM1 = L - 1
DL = 1
                    (L .EQ. N) GO TO 15
(B(L+1,L) .NE. 0.) DL = 2
= L+DL-1
    15
                 (L .EQ. 1) GO TO 30

) 20 J = L,LL

DO 20 I = 1,M

DO 20 IR = 1,LM1

C(I,J) = C(I,J) - C(I,IB)*B(IB,J)
             DO
             CONTINUE
K = 1
KM1 = K
                 1 = K - 1

= 1

(K .EQ. M) GO TO 45

(A(K,K+1) .NE. 0.) DK = 2

= K+DK-1

(K .EQ. 1) GO TO 60

50 I=K,KK

DO 50 J=L,LL

DO 50 JA=1,KM1

C(I,J) = C(I,J) - A(I,JA)*C(JA,J)
                         K - 1
             DK
IF
IF
            C(I,J) = C(I,J) - A(I,JA)*C(CONTINUE)
IF (DL .EQ. 2) GO TO 80
IF (DK .EQ. 2) GO TO 70
I(1,1) = A(K,K) + B(L,L)
IF (T(1,1) .EQ. 0.) RETURN
C(K,L) = C(K,L) / T(1,1)
IF (ABS(C(K,L)) .GE. RMAX) RETURN
GO TO 100
I(1,1) = A(K,K) + B(L,L)
I(1,2) = A(K,KK)
I(2,1) = A(K,KK)
I(2,2) = A(K,KK) + B(L,L)
P(I) = C(K,L)
    70
```

```
P(2) = C(KK,L)

NSYS = 2

CALL SYSSLV

IF (SING) RETURN

C(KK,L) = P(2)

IF (ABS(C(K,L)) .GE. RMAX) RETURN

GO TO 100

80 If (1,2) = B(LL,L)

T(2,1) = B(L,L)

T(2,1) = B(L,L)

P(1) = C(K,L)

P(2) = C(K,L)

P(2) = C(K,L)

NSYS = 2

CALL SYSSLV

IF (SING) RETURN

C(K,L) = P(2)

IF (ABS(C(K,L)) .GE. RMAX) RETURN

GO TO 100

90 I(1,1) = A(K,K) + B(L,L)

T(1,2) = A(K,K)

T(2,2) = A(K,K)

T(2,2) = A(K,K)

T(2,2) = A(K,K) + B(L,L)

T(3,2) = 0

T(3,1) = B(L,L)

T(3,2) = 0

T(4,2) = T(1,3)

T(3,1) = B(L,L)

T(3,2) = O(K,L)

P(1) = C(K,L)

P(2) = C(K,L)

P(3) = C(K,L)

P(4) = C(K,L)

P(5) = C(K,L)

P(5) = C(K,L)

P(1) = C(K,L)

P(2) = C(K,L)

P(3) = C(K,L)

P(4) = C(K,L)

P(5) = C(K,L)

P(1) = C(K,L)

P(2) = C(K,L)

P(3) = C(K,L)

P(4) = C(K,L)

P(5) = C(K,L)

P(6) = RMAX) RETURN

C(K,L) = P(1)

IF (ABS(C(K,L)) .GE. RMAX) RETURN

C(K,L) = P(2)

IF (ABS(C(K,L)) .GE. RMAX) RETURN

C(K,L) = P(3)

IF
                                                                                                                                           P(2) = C(KK \cdot L)
                                                                                                                                                              END
```

```
SUBROUTINE SYSSLY
          SYSSLV IS A FORTRAN IV SUBROUTINE THAT SOLVES THE LINEAR SYSTEM AX = B OF ORDER N LESS THAN 5 BY CROUT REDUCTION FOLLOWED BY BLOCK SUBSTITUTION. THE MATRIX A, THE VECTOR B, AND THE ORDER N ARE CONTAINED IN THE ARRAYS A, B, AND THE VARIABLE N OF THE COMMON BLOCK SLVBLK. THE SOLUTION IS RETURNED IN THE ARRAY B.
           COMMON /SLVBLK/ A(5.5), B(5), N, SING REAL MAX LOGICAL SING SING = .TRUE.

1 NM1 = N - 1 N1 = N + 1
0000
           COMPUTE THE LU FACTORIZATION OF A
                DO 80 K=1,N

KM1 = K-1

IF (K .EQ. 1) GO TO 20

DO 10 I=K,N

DO 10 J=1,KM1

A(I,K) = A(I,K) - A(I,J)*A(J,K)
                       CONTINUE

IF (K .EQ. N) GO TO 100

KP1 = K+1

MAX = ABS(A(K,K))
     20
                      MAX = ABS(A(K,K))

INTR = K

DO 30 I=KP1,N

AA = ABS(A(I,K))

IF (AA .LE. MAX) GO TO 30

MAX = AA

INTR = I

CONTINUE

IF (MAX .EQ. 0.) RETURN

A(N1,K) = INTR

IF (INTR .EQ. K) GO TO 50

DO 40 J=1,N

TEMP = A(INTR,J)

A(K,J) = A(INTR,J)

A(INTR,J) = TEMP
      30
                       A(INTR,J) = TEMP
CONTINUE
DO 80 J=KP1,N
IF (K .EQ. 1) GO TO 70
DO 60 I=1,KM1
      40
                              A(K,J) = A(K,J) - A(K,I)*A(I,J)
      60
70
                 CONTINUE = A(K,J) / A(K,K)
            INTERCHANGE THE COMPONENTS OF B
    100 DO 110 J=1:NM1

INTR = A(N1:J)

IF (INTR .EQ. J) GO TO 110

TEMP = B(J)

B(J) = B(INTR)

B(INTR) = TEMP

110 CONTINUE
            SOLVE SY = B
     200 B(1) = B(1) / A(1,1)

D0 220 I=2,N

IM1 = I-1

D0 210 J=1,IM1

B(I) = B(I) - A(I,J)*B(J)

210 CONTINUE

B(I) = B(I) / A(I,I)

220 CONTINUE
           SOLVE UX=Y
```

C
300 DO 310 II=1.NM1
I = NM1-II+1
II = I+1
DO 310 J=I1.N
B(I) = B(I) - A(I.J)\*B(J)
310 CONTINUE
SING = .FALSE.
RETURN
END

```
SUBROUTINE HOR3 (A, V, N, NLOW, NUP, EPS, ER, FI, TYPE, NA, NV)
C
                      INTEGER N.NA.NLOW.NUP.NV.TYPE(N)
REAL A(NA.N).EI(N).ER(N).EPS.V(NV.N)
                     HOR3 REDUCES THE UPPER HESSENBERG MATRIX A TO QUASITRIANGULAR FORM BY UNITARY SIMILARITY TRANSFORMATIONS. THE EIGENVALUES OF A, WHICH ARE CONTAINED IN THE 1X1 AND 2X2 DIAGONAL BLOCKS OF THE REDUCED MATRIX, ARE ORDERED IN DESCENDING ORDER OF MAGNITUDE ALONG THE DIAGONAL. THE TRANSFORMATIONS ARE ACCUMULATED IN THE ARRAY V. HOR3 REQUIRES THE SUBROUTINES EXCHNG, QRSTEP, AND SPLIT. THE PARAMETERS IN THE CALLING SEQUENCE ARE (STARRED PARAMETERS ARE ALTERED BY THE SUBROUTINE)
AN ARRAY THAT INITIALLY CONTAINS THE N X N UPPER HESSENBERG MATRIX TO BE REDUCED. ON RETURN, A CONTAINS THE REDUCED. QUASITIANGULAR MATRIX.

AN ARRAY THAT CONTAINS A MATRIX INTO WHICH THE REDUCING TRANSFORMATIONS ARE TO BE
                                  *4
                                  * V
                                                                  AN ARRAY THAT CONTAINS A MATRIX INTO WHICH
THE REDUCING TRANSFORMATIONS ARE TO BE
MULTIPLIED.
THE ORDER OF THE MATRICES A AND V.
A(NLOW-1,NLOW) AND A(NUP,NUP+U) ARE
ASSUMED TO BE ZFRO, AND ONLY ROWS NIOW
THROUGH NUP AND COLUMNS NLOW THROUGH
NUP ARE TRANSFORMED, RESULTING IN THE
CALCULATION OF EIGENVALUES NLOW
THROUGH NUP.
A CONVERGENCE CRITERION.
AN ARRAY THAT ON RETURN CONTAINS THE REAL
PARTS OF THE EIGENVALUES.
AND INTEGER ARRAY WHOSE I-TH ENTRY IS
OUT OF THE I-TH FIGENVALUE IS REAL
THE I-TH FIGENVALUE IS COMPLEX
WITH POSITIVE IMAGINARY PART.
IF THE I-TH FIGENVALUE IS COMPLEX
WITH NEGATIVE IMAGINARY PART.
THE I-TH FIGENVALUE WAS NOT
CALCULATED SUCCESSFULLY.
THE FIRST DIMENSION OF THE ARRAY A.
THE FIRST DIMENSION OF THE ARRAY A.
                                     NLOW
                                     EPS
                                  *ER
                                  *EI
                                  *TYPE
                                     NA
                      INTERNAL VARIABLES
                      INTEGER I.IT.L.MU.NL.NU
REAL E1.E2.P.G.R.S.T.W.X.Y.Z
LUGICAL FAIL
                      INITIALIZE.
           DU 10 I=NLOW.NUP
TYPE(I) = -1
                       T = 0.
CCC
                      MAIN LOOP. FIND AND ORDER EIGENVALUES.
       100 IF (NU .LT. NLOW) GO TO 500
0000
                             OR LOOP.
                                                                   FIND NEGLIGABLE ELFMENTS AND PERFORM
                              CONTINUE
       110
                                      SEARCH BACK FOR NEGLIGABLE ELEMENTS.
                                     CONTINUE
       120
```

```
IF(L .EQ. NLOW) GO TO 130
                        IF(ABS(A(L,L-1)) .LT. EPS*(ABS(A(L-1,L-1))+ABS(A(L,L))))
GO TO 130
           1
                    GO TO 120
                    CONTINUE
      130
  0000
                    TEST TO SEE IF AN EIGENVALUE OR A 2X2 BLOCK HAS BEEN FOUND.
                    X = A(NU,NU)

IF(L .EQ. NU) GO TO 300

Y = A(NU-1,NU-1)

W = A(NU,NU-1)*A(NU-1,NU)

IF(L .EQ. NU-1) GO TO 200
                     TEST ITERATION COUNT. IF IT IS 30 QUIT. IT IS 10 OR 20 SET UP AN AD-HOC SHIFT.
                     IF(IT .EQ. 60) GO TO 500
IF(IT.NE.40 .AND. IT.NE.50) GO TO 150
  CCC
                        AD-HOC SHIFT.
                        T = T + X

DO 140 I=NLOW, NU

A(I,I) = A(I,I) - X

CONTINUE

S = ABS(A(NU,NU-1)) + ABS(A(NU-1,NU-2))

X = 0.75±S
      140
                        X = 0.75*S
Y = X
W = -0.437
                              -0.4375*5**2
                     CONTINUE + 1
      150
  0000
                    LOOK FOR TWO COMSECUTIVE SMALL SUB-DIAGONAL ELEMENTS.
                     CONTINUE NU-S
      160
                              NOE

A(NL,NL)

X - Z

Y - Z

(R*S-W)/A(NL+1,NL) + A(NL,NL+1)

A(NL+1,NL+1) - Z - R - S

A(NL+2,NL+1)

AHS(P) + ABS(Q) + ABS(R)

P/S

Q/S
                    P = P/S
Q = Q/S
R = R/S
IF(NL .EQ. L) GO TO 170
IF(ABS(A(NL,NL-1))*(ABS(Q)+ABS(R))
EPS*AHS(P)*(ABS(A(NL-1,NL-1))+AF
GO TO 170
NL = NL-1
GO TO 160
CONTINUE
                                                                    (Q)+ABS(R)) .LF.
-1,NL-1))+ABS(7)+ABS(A(NL+1,NL+1))))
      170
  CCC
                     PERFORM A OR STEP BETWEEN NL AND NU.
                 GO TO 110
                 2X2 BLOCK FOUND.
                 IF(NU .NE. NLOW+1) A(NU-1.NU-2) = 0.
A(NU,NU) = A(NU.NU) + T
A(NU-1.NU-1) = A(NU-1.NU-1) + T
      200
                 TYPE(NU) = 0
TYPE(NU-1) = 0
MU = NU
C 210
                 LOOP TO POSITION 2X2 BLOCK.
                 CONTINUE
                    NL = MU-1
```

```
CALL EXCHNG(A.V.N.MU.1.2.EPS.FAIL.NA.NV)
                         IF(.NOT. FAIL) GO TO 325
TYPE(MU) = -1
                             TYPE (MU+1) = -1
                        GO TO 500
CONTINUE
MU = MU+2
GO TO 340
    325
                         CONTINUE
    330
                             THE NEXT BLOCK IS 1X1.
                    IF (ABS (A (MU, MU)) .GE. ABS (A (MU+1, MU+1)))
GO TO 350

CALL EXCHNG (A, V, N, MU, 1, 1, EPS, FAIL, NA, NV)
MU = MU+1
CONTINUE
MU = NL
NL = 0
(MU, NE = 0) GO TO 310
          1
    340
    350
                 IF (MU .NE. 0) GO TO 310
CCC
                GO HACK AND GET THE NEXT EIGENVALUE.
    400
                CONTINUE
            60 TO 100
CCC
            ALL THE FIGHVALUES HAVE BEEN FOUND AND ORDERED.
    500 IF(NU .LT. NLOW) GO TO 507

DO 503 1=1.NU

A(I,I) = A(I,I) + T

503 CONTINUE
   503 CONTINUE
507 CONTINUE
NU = NUP
510 CONTINUE
IF (TYPE (NU) •NE• -1) GO TO 515
NU = NU-1
                NU = NU-1

GO TO 540

CONTINUE

IF(NU .EQ. NLOW) GO TO 520

IF(A(NU,NU-1) .EQ. 0.) GO TO 520
                     2X2 BLOCK .
                     CALL SPLIT (A, V.N.NU-1, E1, E2, NA, NV)
IF (A(NU, NU-1) . E0. 0.) 60 TO 520
                     FR(NU) = E1

ER(NU) = E2

ER(NU-1) = ER(NU)

EI(NU) = -EI(NU-1)

TYPE(NU-1) = 1

TYPE(NU) = 2
                NU = NU-2
GO TO 530
CONTINUE
    520
           ER(NU) = A(NU,NU)
EI(NU) = 0.
NU = NU-1
CONTINUE
CONTINUE
IF(NU .GE. NLOW) GO TO 510
RETURN
FND
```

```
C
                ATTEMPT TO SPLIT THE BLOCK INTO TWO REAL
CCC
               EIGENVALUES.
                CALL SPLIT (A.V.N.NL.E1, E2, NA.NV)
CCCC
               IF THE SPLIT WAS SUCCESSFUL, GO AND ORDER THE
                IF(A(MU.MU-1) .EQ. 0.) GO TO 310
                TEST TO SEE IF THE BLOCK IS PROPERLY POSITIONED.
                IF(MU .EQ. NUP) GO TO 400
IF(MU .EQ. NUP-1) GO TO 220
IF(A(MU+2,MU+1) .EQ. U.) GO TO 220
                   THE NEXT BLOCK IS 2X2.
                   IF (A (MU-1, MU-1) *A (MU, MU) -A (MU-1, MU) *A (MU, MU-1)
                       • GE • A(MU+1, MU+1) *A(MU+2, MU+2) -A(MU+1, MU+2) *

A(MU+2, MU+1))

GO TO 400
        123
                  CALL EXCHNG(A,V,N,NL,2,2,EPS,FAIL,NA,NV)
IF(.NOT. FAIL) GO TO 215
TYPE(NL) = -1
TYPE(NL+1) = -1
TYPE(NL+2) = -1
TYPE(NL+3) = -1
GO TO 500
CONTINUE
               CONTINUE
MU = MU+2
GO TO 230
CONTINUE
   215
   220
CCC
                   THE NEXT BLOCK IS 1X1.
                   IF (A(MU-1, MU-1) *A(MU, MU) -A(MU-1, MU) *A(MU, MU-1)
                        .GF . A(MU+1, MU+1) **2)
GO TO 400
        2
                        GO
                   CALL EXCHNG(A, V, N, NL, 2, 1, EPS, FAIL, NA, NV)

IF(.NOT. FAIL) GO TO 225

TYPE(NL) = -1

TYPE(NL+1) = -1

TYPE(NL+2) = -1

GO TO 500
                  CONTINUE
MU = MU+1
   225
            GO TO 210
   230
CCC
             SINGLE EIGENVALUE FOUND.
   300
             A(NU,NU) = A(NU,NU) + T
            IF(NU .NE. NLOW) A(NU,NU-1) = 0.

TYPE(NU) = 0

MU = NU
CCC
            LOOP TO POSITION ONE OR TWO REAL EIGENVALUES.
             CONTINUE
   310
                POSITION THE EIGENVALUE LOCATED AT A(NL.NL).
Č
   320
                CONTINUE
                   IF(MU .EQ. NUP) GO TO 350
IF(MU .EQ. NUP-1) GO TO 330
IF(A(MU+2,MU+1) .EQ. 0.) GO TO 330
                      THE NEXT BLOCK IS 2X2.
                      IF(A(MU, MU) **2 .GE.
A(MU+1, MU+1) *A(MU+2, MU+2) -A(MU+1, MU+2) *A(MU+2, MU+1))
GO TO 400
        ž
```

```
SUBROUTINE EXCHNG (A.V.N.L.B1.B2.EPS.FAIL.NA.NV)
C
                       INTEGER B1.B2.L.NA.NV
REAL A(NA.N).EPS,V(NV.N)
LOGICAL FAIL
GIVEN THE UPPER HESSENBERG MATRIX A WITH CONSECUTIVE B1XB1 AND B2XB2 DIAGONAL BLOCKS (B1,B2 .LE. 2) STARTING AT A(L,L), EXCHNG PRODUCES A UNITARY SIMILARITY TRANSFORMATION THAT EXCHANGES THE PLOCKS ALONG WITH THEIR EIGENVALUES. THE TRANSFORMATION IS ACCUMULATED IN V. EXCHNG REQUIRES THE SUBROUTINE GRSTEP. THE PARAMETERS IN THE CALLING SEQUENCE ARE (STARRED PARAMETERS ARE ALTERED BY THE SUBROUTINE)
                                                                    THE MATRIX WHOSE BLOCKS ARE TO RE INTERCHANGED.
THE ARRAY INTO WHICH THE TRANSFORMATIONS ARE TO RE ACCUMULATED.
THE ORDER OF THE MATRIX A.
THE POSITION OF THE BLOCKS.
THE SIZE OF THE FIRST BLOCK.
THE SIZE OF THE SECOND BLOCK.
A CONVERGENCE CRITERION.
A LOGICAL VARIABLE WHICH IS FALSE ON A NORMAL RETURN. IF THIRTY ITERATIONS WERE PERFORMED WITHOUT CONVERGENCE, FAIL IS SET TO TRUE AND THE ELFMENT
A(L+B2,L+B2-1) CANNOT BE ASSUMED ZERO.
THE FIRST DIMENSION OF THE ARRAY A.
THE FIRST DIMENSION OF THE ARRAY V.
                                   * 4
                                   * V
                                       N
                                  B1
B2
EPS
*FAIL
                                      NA
                       INTERNAL VARIABLES.
                       INTEGER I, IT, J, L1, M
REAL P, Q, R, S, W, X, Y, Z
C
                       FAIL = .FALSE.
IF(B1 .EQ. 2) GO TO 40
IF(B2 .EQ. 2) GO TO 10
                                       INTERCHANGE 1X1 AND 1X1 BLOCKS.
                                            1 = L+1
= A(L+1,L+1) - A(L,L)
= A(L,L+1)
= AMAX1(P,Q)
                                       L1
                                       GP
                                      IF (R .E0. 0.) RETURN
P = P/R
0 = 0/R
R = SQRT(P**2 + Q**2)
P = P/R
                                      P = P/R

Q = Q/R

DO 3 J=L,N

S = P*A(L,J) + Q*A(L+1,J)

A(L+1,J) = P*A(L+1,J) - Q*A(L,J)

A(L,J) = S

CONTINUE

DO 5 I=1,L1

S = P*A(I,L) + Q*A(I,L+1)

A(I,L) = P*A(I,L+1) - Q*A(I,L)

A(I,L) = S
                3
                                      CONTINUÉ

DO 7 I=1,N

S = P*V(I,L) + Q*V(I,L+1)

V(I,L+1) = P*V(I,L+1) - Q*V(I,L)

V(I,L) = S
                               CONTINUE
A(L+1,L) = 0.
RETURN
CONTINUE
               7
            10
                                       INTERCHANGE 1X1 AND 2X2 BLOCKS.
```

```
X = A(L,L,

P = 1.

Q = 1.

R = 1.

CALL QRSTEP(A,V,P,Q,R,L,L+2,N,NA,NV)

IT = Q

IT = IT+1

IF(IT ,LE. 60) GO TO 30

FAIL = .TRUE.

RETURN

CONTINUE X
                  X = A(L,L)
     20
                     RETURN
CONTINUE
P = A(L,L) - X
Q = A(L+1,L)
R = 0
CALL GRSTEP(A,V,P,Q,R,L,L+2,N,NA,NV)
IF(ABS(A(L+2,L+1)) .GT.
EPS*(AHS(A(L+1,L+1))+ABS(A(L+2,L+2))))
     30
     GO TO 20
A(L+2,L+1) = 0.

RETURN

40 CONTINUE
000
               INTERCHANGE 2X2 AND B2XB2 BLOCKS.
               M = L+2

IF(B2 .EQ. 2) M = M+1

X = A(L+1,L+1)

Y = A(L,L)

W = A(L+1,L)*A(L,L+1)
          50
      60
```

```
C
                     INTEGER L,N,NA,NV
REAL A(NA,N),V(NV,N)
\sigma
                    GIVEN THE UPPER HESSENBERG MATRIX A WITH A 2X2 BLOCK STARTING AT A(L,L), SPLIT DETERMINES IF THE CURRESPONDING EIGENVALUES ARE REAL OR COMPLEX. IF THEY ARE REAL, A ROTATION IS DETERMINED THAT REDUCES THE BLOCK TO UPPER TRIANGULAR FORM WITH THE FIGENVALUE OF LARGEST ABSOLUTE VALUE APPEARING FIRST. THE HOTATION IS ACCUMULATED IN V. THE EIGENVALUES (REAL OR COMPLEX) ARE RETURNED IN E1 AND E2. THE PARAMETERS IN THE CALLING SEQUENCE ARE (STARRED PARAMETERS ARE ALTERED BY THE SUBROUTINE)
                                                                THE UPPER HESSENVERG MATRIX WHOSE 2X2
BLOCK IS TO BE SPLIT.
THE ARRAY IN WHICH THE SPLITTING TRANS-
FORMATION IS TO BE ACCUMULATED.
THE ORDER OF THE MATRIX A.
THE POSITION OF THE 2X2 BLOCK.
ON RETURN IF THE EIGENVALUES ARE COMPLEX
E1 CONTAINS THEIR COMMON REAL PART AND
E2 CONTAINS THE POSITIVE IMAGINARY PART.
IF THE EIGENVALUES ARE REAL, E1 CONTAINS
THE ONE LARGEST IN ABSOLUTE VALUE AND F2
CONTAINS THE OTHER ONE.
THE FIRST DIMENSION OF THE ARRAY A.
THE FIRST DIMENSION OF THE ARRAY V.
                                 * A
                                 * V
                                    N
                                 *E1
                                    NV
                      INTERNAL VARIABLES
                     INTEGER I.J.L1
REAL P.Q.R.T.U.W.X.Y.Z
C
                           = A(L+1,L+1)
= A(L,L)
= A(L,L+1)*A(L+1,L)
                           = (Y-X)/2.
= P**2 + W
                      IF (Q .GE. 0.) GO TO 5
                             COMPLEX EIGENVALUE.
              E1 = P + X
E2 = SQRT(-Q)
RETURN
5 CONTINUE
                                                                                                         SET UP TRANSFORMATION.
                      TWO REAL EIGENVALUES.
         Z = SQRT(Q)

IF (P .LT. 0.) GO TO 10

Z = P + Z

GO TO 20

10 CONTINUE

Z = P - Z

20 CONTINUE

LF (Z .FQ. 0.) GO TO 30
                      IF(Z .EQ. 0.) GO TO 30
          30 CONTINUE
                    CONTINUE

R = 0.

CONTINUE

IF (ABS(X+Z) .GE. ABS(X+R)) Z = R

Y = Y - X - Z

X = -Z

T = A(L,L+1)

U = A(L+1,L)

U = A(L+1,L)
                      U = A(L+1,L)
IF(ABS(Y)+ABS(U) .LE. ABS(T)+ABS(X)) GO TO 60
                             GO TO 70
```

SUBROUTINE SPLIT (A, V, N, L, E1, E2, NA, NV)

```
SUBROUTINE GRSTEP(A, V, P, Q, R, NL, NU, N, NA, NV)
C
                      INTEGER N.NA.NL.NU.NV
REAL A(NA.N).P.G.R.V(NV.N)
QRSTEP PERFORMS ONE IMPLICIT OR STEP ON THE UPPER HESSENBERG MATRIX A. THE SHIFT IS DETERMINED BY THE NUMBERS P.G. AND R. AND THE STEP IS APPLIED TO HOWS AND COLUMNS NL THROUGH NU. THE TRANSFORMATIONS ARE ACCUMULATED IN V. THE PARAMETERS IN THE CALLING SEQUENCE ARE (STARRED APRAMETERS ARE ALTERED BY THE SUBROUTINE)
                                                                     THE UPPER HESSENBERG MATRIX ON WHICH THE OR STEP IS TO HE PERFORMED. THE ARRAY IN WHICH THE TRANSFORMATIONS ARE TO HE ACCUMULATED PARAMETERS THAT DETERMINE THE SHIFT.
                                  * 4
                                   *٧
                                   *P
                                   *0
*R
                                      NL
                                                                     THE LOWER LIMIT OF THE STEP.
THE UPPER LIMIT OF THE STEP.
THE ORDER OF THE MATRIX A.
THE FIRST DIMENSION OF THE ARRAY A.
THE FIRST DIMENSION OF THE ARRAY V.
                                      N
                       INTERNAL VARIABLES.
                        INTEGER I.J.K.NL2.NL3.NUM1
                       REAL S.X.Y.Z
LUGICAL LAST
C
           NL2 = NL+2

DO 10 I=NL2,NU

A(I,I-2) = 0.

10 CONTINUE
           10 CONTINUE

IF (NL2 .EQ. NU) GO TO 30

NL3 = NL+3

DO 20 I=NL3.NU

A(I,I-3) = 0.

20 CONTINUE

30 CONTINUE

NUM1 = NU-1

DO 130 K=NL.NUM1
CCC
                            DETERMINE THE TRANSFORMATION.

LAST = K .EQ. NUM1

IF (K .EQ. NL) GO TO 40

P = A(K,K-1)

Q = A(K+1,K-1)

R = 0.

IF (.NOT.LAST) R = A(K+2,K-1)

X = ABS(P) + ABS(Q) + APS(R)

IF (X .EQ. 0.) GO TO 130

P = P/X

Q = Q/X

R = R/X

CONTINUE

S = SQRT(P**2 + Q**2 + R**2)

IF (K .EQ. NL) GO TO 50

A(K,K-1) = -S*X

GO TO 60

CONTINUE

IF (NL .NE. 1) A(K,K-1) = -A(K,K-1)

CONTINUE

P = P + S

X = P/S

Y = Q/S

Z = R/S

Q = Q/P

R = R/P
                               DETERMINE THE TRANSFORMATION.
           40
           50
           60
C
```

```
PREMULTIPLY.
C
                  DO BU J=K,N
P = A(K,J) + Q*A(K+1,J)

IF(LAST) GO TO 70
P = P + R*A(K+2,J)
A(K+2,J) = A(K+2,J) - P*Z

CONTINUE
A(K+1,J) = A(K+1,J) - P*Y
A(K,J) = A(K,J) - P*X

CONTINUE
C
       70
       80
                   POSTMULTIPLY.
                   J = MINO(K+3,NU)

DO 100 I=1,J

P = X*A(I,K) + Y*A(I,K+1)

IF(LAST) GO TO 90

P = P + Z*A(I,K+2)

A(I,K+2) = A(I,K+2) - P*R

CONTINUE

A(I,K+1) = A(I,K+1) - P*Q

A(I,K) = A(I,K) - P
        90
                    CONTINUE

ACCUMULATE THE TRANSFORMATION IN V.
      100
 CCC
                    UO 120 I=1,N

P = X*V(I,K) + Y*V(I,K+1)

IF(LAST) GO TO 110

P = P + Z*V(I,K+2)

V(I,K+2) = V(I,K+2) - P*R
                    CONTINUE

V(I,K+1) = V(I,K+1) - P*Q

V(I,K) = V(I,K) - P

CONTINUE
       110
       120
       130 CONTINUE
                RETURN
END
```

```
SUBROUTINE ORTHES(NM,N,LOW,IGH,A,ORT)
INTEGER I,J,M,N,II,JJ,LA,MP,NM,IGH,KP1,LOW
REAL A(NM,N),ORT(IGH)
REAL F,G,H,SCALE
LA = IGH = 1
KP1 = LOW + 1
IF(LA .LT. KP1) GO TO 200

DO 180 M=KP1,LA

H = 0.
SCALE = 0.
ORT(M) = 0.
SCALE = SCALE + ARS(A(I,M-1))
IF(SCALE .EQ. 0.) GO TO 180

MP = M + IGH
DO 100 II=M,IGH
I = MP = II
ORT(I) = A(I,M-1)/SCALE
H = H + ORT(I)*ORT(I)
CONTINUE
G = -SIGN(SGRT(H),ORT(M))
H = H - ORT(M)*G
ORT(M) = ORT(M) - G
DO 130 J=M,N
F = 0.
DO 110 II=M,IGH
I = MP - II
F = F + ORT(I)*A(I,J)
CONTINUE
IF (H .EQ.0.) GO TO 162
F = F/H
DO 120 I=M,IGH
A(I,J) = A(I,J) - F*ORT(I)
CONTINUE
CONTINUE
CONTINUE
CONTINUE
DO 160 I=1,IGH
         90
  100
120
130
```

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19. KEY WORDS (Continue on reverse side if necessary and identity by block of reducing subspaces eigenvalues block diagonalization eigenvectors.  ABSTRACT (Continue on reverse side if necessary and identity by block of the diagonal form by a real similarity transformation formation corresponding to a block span a red block is the representation of A in that su bas is. The algorithm attempts to control the tion matrices, so that the reducing subspaces	g a real matrix A to block tion. The columns of the transucing subspace of A, and the bspace with respect to the condition of the transforma-